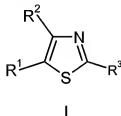


**In the Claims:**

The current status of all claims is listed below and supersedes all previous lists of claims.

Please cancel claims 7-11 without prejudice to their presentation in another application, and amend claims 1-3 as follows:

1. (currently amended) A compound of formula (I)



wherein

$R^1$  and  $R^2$  are ~~independently selected from phenyl, thienyl, and pyridyl~~, each of which is independently optionally substituted with one, two or three Z groups;

Z is selected from a  $C_{1-6}$ alkyl group, a  $C_{1-6}$ alkoxy group, hydroxy, halo, trifluoromethyl, trifluoromethylthio, trifluoromethoxy, trifluoromethylsulphonyl, nitro, amino, mono or di  $C_{1-3}$ alkylamino, mono or di  $C_{1-3}$ alkylamido,  $C_{1-3}$ alkylsulphonyl,  $C_{1-3}$ alkoxycarbonyl, carboxy, cyano, carbamoyl, mono or di  $C_{1-3}$ alkyl carbamoyl, sulphamoyl, acetyl,  $-O-CH_2-CH_2-O-$  attached at two adjacent carbons, and phenyl, optionally substituted with one or more of the following: a  $C_{1-6}$ alkyl group, trifluoromethyl, a  $C_{1-6}$ alkoxy group, trifluoromethoxy, halo, or  $-O-CH_2-CH_2-O-$  attached at two adjacent carbons;

$R^3$  is  $-X-Y-NR^4R^5$ ;

$R^4$  and  $R^5$  are independently selected from:

a  $C_{1-6}$ alkyl group, optionally substituted with a  $C_{1-6}$ alkoxy group or trifluoromethoxy;

an (amino) $C_{1-4}$ alkyl group, wherein the amino is optionally substituted by one or more  $C_{1-3}$ alkyl groups;

a non-aromatic  $C_{2-15}$ carbocyclic group, optionally substituted with a  $C_{1-3}$ alkoxy $C_{1-3}$ alkyl group;

a-(C<sub>2-12</sub>cycloalkyl)C<sub>1-3</sub>alkyl-group;

a-(CH<sub>2</sub>)<sub>r</sub>(phenyl)<sub>s</sub> group, wherein r is 0, 1, 2, 3 or 4, and wherein s is 1 when r is 0, otherwise s is 1 or 2, and wherein the phenyl groups are optionally independently substituted with one, two or three Z groups;

naphthyl;

anthracenyl;

a saturated 5- to 8-membered heterocyclic piperidine group containing one ~~nitrogen and~~ optionally containing one of the following: oxygen, sulphur or an additional nitrogen, wherein the ~~heterocyclic~~ piperidine group is optionally substituted by one or more C<sub>1-3</sub>alkyl groups or benzyl;

1-adamantylmethyl; and

a -(CH<sub>2</sub>)<sub>t</sub>Het group, wherein t is 0, 1, 2, 3 or 4, and the alkylene chain is optionally substituted by one or more C<sub>1-3</sub>alkyl groups, and wherein Het is ~~an aromatic heterocycle optionally substituted by one, two or three groups selected from a C<sub>1-6</sub>alkyl group; a C<sub>1-6</sub>alkoxy group, trifluoromethoxy or halo or~~ Het is a saturated 5- to 8-membered heterocyclic a piperidine group ~~containing one nitrogen and~~ optionally containing one of the following: oxygen, sulphur or an additional nitrogen; wherein the ~~heterocyclic~~ piperidine group is optionally substituted by one or more C<sub>1-3</sub>alkyl groups, hydroxy or benzyl; and

wherein R<sup>4</sup> may be H; and

wherein R<sup>4</sup> and R<sup>5</sup> taken together with the nitrogen atom to which they are attached form a saturated 5- to 8-membered heterocyclic piperidine group ~~containing one nitrogen and~~ optionally containing one of the following: oxygen, sulphur or an additional nitrogen; wherein the ~~heterocyclic~~ piperidine group is optionally substituted with one or more C<sub>1-3</sub>alkyl groups, hydroxy or benzyl;

X is CO or SO<sub>2</sub>; and

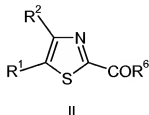
Y is absent or represents NH optionally substituted by a C<sub>1-3</sub>alkyl group;

or a pharmaceutically acceptable salt, ~~prodrug or solvate~~ thereof;

with the proviso that R<sup>1</sup> and R<sup>2</sup> are not both 4-methoxyphenyl and the proviso that when

$R^1$  is phenyl and  $R^2$  represents phenyl or 4-fluorophenyl, X is CO and Y is absent then the group  $NR^4R^5$  is not methyl-[2-[1-(phenylmethyl)-4-piperidinyl]ethyl]amino, methylpiperazino, 2-[1-methyl-4-piperidinyl]ethylamino, or [2-[1-(phenylmethyl)-4-piperidinyl]ethyl]amino.

2. (currently amended) A compound of formula I as represented by formula (II)



wherein

$R^1$  is phenyl, optionally substituted by one or more of the following: a  $C_{1-6}$ alkyl group, trifluoromethyl, a  $C_{1-6}$ alkoxy group, trifluoromethoxy, halo, or -O-CH<sub>2</sub>-CH<sub>2</sub>-O- attached at two adjacent carbons;

$R^2$  is phenyl, optionally substituted by one or more of the following: a  $C_{1-6}$ alkyl group, trifluoromethyl, a  $C_{1-6}$ alkoxy group, trifluoromethoxy, halo, or -O-CH<sub>2</sub>-CH<sub>2</sub>-O- attached at two adjacent carbons; and

$R^6$  is selected from 1-piperidinylamino, a  ~~$C_{1-6}$ cycloalkylamino group, optionally substituted by  $C_{1-3}$ alkoxy  $C_{1-3}$ alkyl, pyridylamino, wherein the pyridyl ring is optionally substituted by one or more of the following: a  $C_{1-6}$ alkyl group; a  $C_{1-6}$ alkoxy group or trifluoromethoxy; a  $C_{1-6}$ alkylamino group, wherein the alkyl chain is optionally substituted by one or more of the following: a  $C_{1-6}$ alkoxy group, trifluoromethoxy or morpholine;~~  
 or a pharmaceutically acceptable salt ~~prodrug or solvate thereof;~~  
 with the proviso that when  $R^1$  is 4-methoxyphenyl and  $R^2$  is 4-methoxyphenyl, then  $R^6$  is not 2-(morpholine)ethyl.

3. (currently amended) A compound selected from:

4-(4-chlorophenyl)-5-(2,4-dichlorophenyl)thiazole-2-carboxylic acid cyclohexylamide;  
 5-(4-chlorophenyl)-4-(2,4-dichlorophenyl)thiazole-2-carboxylic acid cyclohexylamide;  
 4-(4-chlorophenyl)-5-(2,4-dichlorophenyl)thiazole-2-carboxylic acid piperidin-1-

ylamide;

5-(4-chlorophenyl)-4-(2,4-dichlorophenyl)thiazole-2-carboxylic acid piperidin-1-

ylamide;

4-(4-bromophenyl)-5-phenylthiazole-2-carboxylic acid cyclohexylamide;

4-(4-bromophenyl)-5-phenylthiazole-2-carboxylic acid piperidin-1-ylamide;

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid cyclohexylamide;

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid piperidin-1-ylamide;

4-(4-methoxyphenyl)-5-phenylthiazole-2-carboxylic acid cyclohexylamide;

4,5-bis-(4-methoxyphenyl)thiazole-2-carboxylic acid cyclohexylamide;

4,5-bis-(4-methoxyphenyl)thiazole-2-carboxylic acid piperidin-1-ylamide;

5-(7-bromo-2,3-dihydrobenzo[1,4]dioxin-6-yl)-4-phenylthiazole-2-carboxylic acid piperidin-1-ylamide;

4-(7-bromo-2,3-dihydrobenzo[1,4]dioxin-6-yl)-5-phenylthiazole-2-carboxylic acid

piperidin-1-ylamide;

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid (2-methoxymethylcyclopentyl)-

amide;

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid pyridin-4-ylamide;

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid (2-ethoxyethyl)amide; and

4,5-bis-(4-chlorophenyl)thiazole-2-carboxylic acid (2-morpholin-4-yl-ethyl)amide

and where applicable, optical isomers, tautomers, stereoisomers and racemates thereof as well as pharmaceutically acceptable salts and solvates thereof.

4. (canceled).

5. (previously presented) A pharmaceutical formulation comprising a compound of any one of claims 1 to 3 and a pharmaceutically acceptable adjuvant, diluent or carrier.

6-11. (canceled).

## REMARKS